



# Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 01:51 PM GMT

PDB ID : 3PWP  
Title : The complex between TCR A6 and human Class I MHC HLA-A2 with the bound HuD peptide  
Authors : Borbulevych, O.Y.; Baker, B.M.  
Deposited on : 2010-12-08  
Resolution : 2.69 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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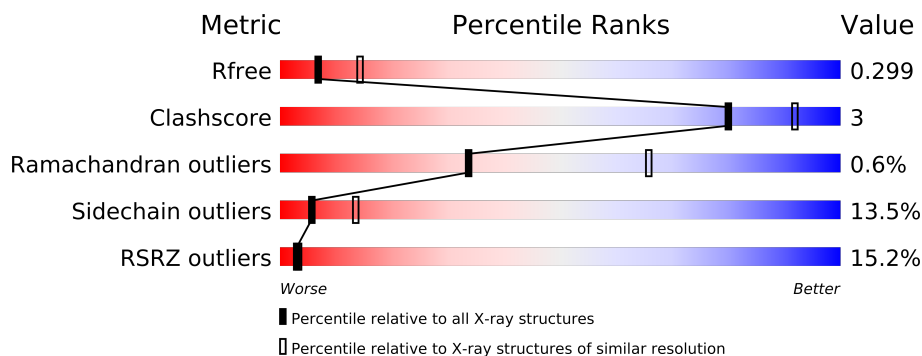
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.69 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	66092	1557 (2.70-2.70)
Clashscore	79885	1939 (2.70-2.70)
Ramachandran outliers	78287	1905 (2.70-2.70)
Sidechain outliers	78261	1905 (2.70-2.70)
RSRZ outliers	66119	1559 (2.70-2.70)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	275	
2	B	100	
3	C	9	
4	D	200	
5	E	245	

## 2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 6757 atoms, of which 0 are hydrogen and 0 are deuterium.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called HLA class I histocompatibility antigen, A-2 alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	275	Total	C	N	O	S	0	0	0
			2247	1403	409	426	9			

- Molecule 2 is a protein called Beta-2-microglobulin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	100	Total	C	N	O	S	0	1	0
			843	537	142	160	4			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	0	MET	-	INITIATING METHIONINE	UNP P61769

- Molecule 3 is a protein called HuD peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	9	Total	C	N	O	0	0	0
			75	52	10	13			

- Molecule 4 is a protein called A6 TCR alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	200	Total	C	N	O	S	0	0	0
			1552	965	255	325	7			

- Molecule 5 is a protein called A6 TCR beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	245	Total	C	N	O	S	0	0	0
			1927	1209	338	372	8			

- Molecule 6 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			6	3	3		
6	A	1	Total	C	O	0	0
			6	3	3		
6	A	1	Total	C	O	0	0
			6	3	3		
6	A	1	Total	C	O	0	0
			6	3	3		
6	E	1	Total	C	O	0	0
			6	3	3		
6	E	1	Total	C	O	0	0
			6	3	3		

- Molecule 7 is SULFATE ION (three-letter code: SO4) (formula:  $O_4S$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	B	1	Total	O	S	0	0
			5	4	1		

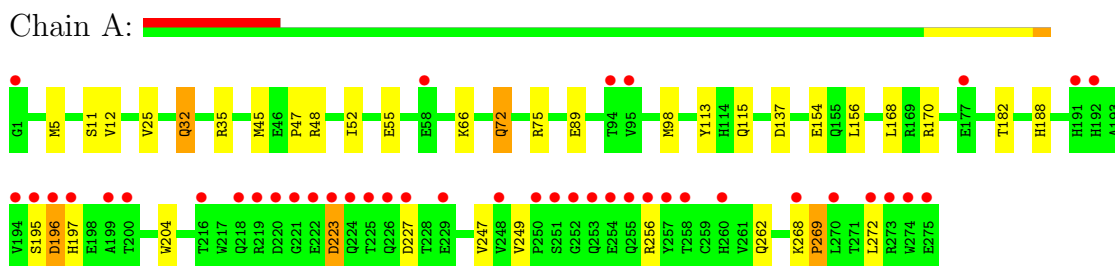
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	19	Total	O	0	0
			19	19		
8	B	16	Total	O	0	0
			16	16		
8	D	7	Total	O	0	0
			7	7		
8	E	24	Total	O	0	0
			24	24		

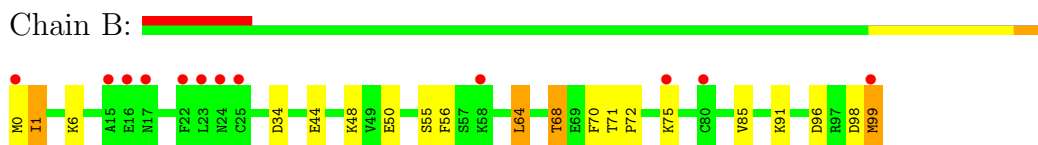
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: HLA class I histocompatibility antigen, A-2 alpha chain



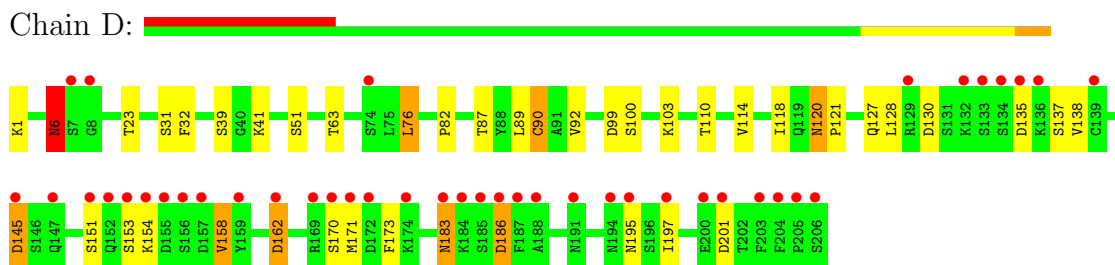
- Molecule 2: Beta-2-microglobulin



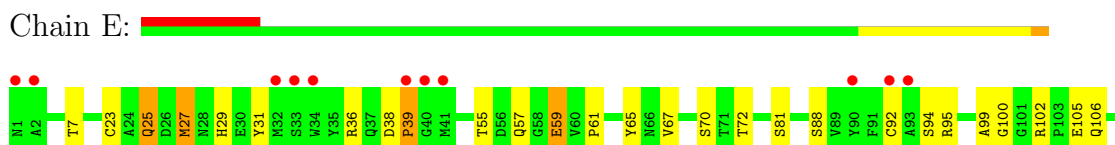
- Molecule 3: HuD peptide

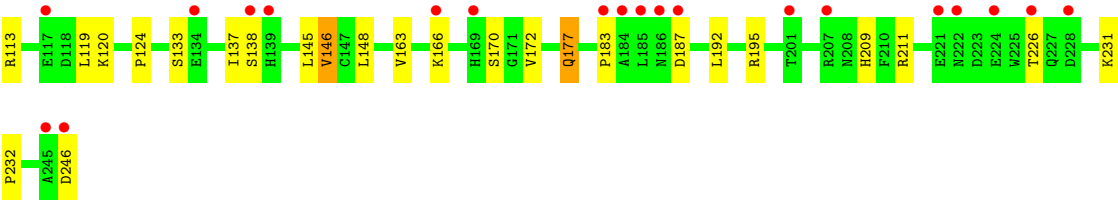


- Molecule 4: A6 TCR alpha chain



- Molecule 5: A6 TCR beta chain





## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	224.01Å 49.06Å 93.71Å 90.00° 90.07° 90.00°	Depositor
Resolution (Å)	20.00 – 2.69 19.83 – 2.69	Depositor EDS
% Data completeness (in resolution range)	94.2 (20.00-2.69) 94.2 (19.83-2.69)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.88 (at 2.71Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.198 , 0.259 0.254 , 0.299	Depositor DCC
$R_{free}$ test set	1370 reflections (5.07%)	DCC
Wilson B-factor (Å <sup>2</sup> )	58.5	Xtriage
Anisotropy	0.207	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 34.5	EDS
Estimated twinning fraction	0.019 for -h,-k,l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 27035 reflections	Xtriage
$F_o, F_c$ correlation	0.90	EDS
Total number of atoms	6757	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.69% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.



## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, SO4

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	A	0.66	0/2312	0.80	1/3137 (0.0%)
2	B	0.65	0/869	0.82	0/1174
3	C	0.76	0/77	0.67	0/102
4	D	0.61	1/1585 (0.1%)	0.83	3/2150 (0.1%)
5	E	0.65	1/1980 (0.1%)	0.84	2/2699 (0.1%)
All	All	0.65	2/6823 (0.0%)	0.82	6/9262 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
5	E	0	2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	90	CYS	CB-SG	-5.51	1.72	1.81
5	E	92	CYS	CB-SG	-5.12	1.73	1.81

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	89	LEU	CA-CB-CG	7.14	131.72	115.30
5	E	23	CYS	CA-CB-SG	-5.40	104.27	114.00
4	D	99	ASP	CB-CG-OD1	5.33	123.10	118.30
1	A	137	ASP	CB-CG-OD1	5.27	123.05	118.30
4	D	76	LEU	CA-CB-CG	5.11	127.04	115.30
5	E	100	GLY	N-CA-C	5.08	125.81	113.10

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
5	E	25	GLN	Peptide
5	E	99	ALA	Peptide

## 5.2 Close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2247	0	2096	13	0
2	B	843	0	811	6	0
3	C	75	0	72	0	0
4	D	1552	0	1461	13	0
5	E	1927	0	1830	14	0
6	A	30	0	40	0	0
6	E	12	0	16	1	0
7	B	5	0	0	0	0
8	A	19	0	0	0	0
8	B	16	0	0	0	0
8	D	7	0	0	0	0
8	E	24	0	0	0	0
All	All	6757	0	6326	44	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 3.

All (44) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
4:D:162:ASP:OD2	4:D:162:ASP:N	2.30	0.63
5:E:95:ARG:HG2	5:E:106:GLN:HB2	1.84	0.59
1:A:25:VAL:HG13	1:A:32:GLN:HE21	1.68	0.57
1:A:188:HIS:HB3	1:A:204:TRP:HB2	1.88	0.55
4:D:118:ILE:HD11	4:D:145:ASP:HA	1.90	0.53
5:E:124:PRO:HD3	5:E:232:PRO:HB3	1.93	0.50
4:D:183:ASN:N	4:D:183:ASN:OD1	2.44	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:55:GLU:OE1	1:A:170:ARG:NH2	2.45	0.49
1:A:5:MET:HB2	1:A:168:LEU:HD13	1.94	0.49
5:E:57:GLN:HB2	5:E:61:PRO:HB3	1.95	0.48
1:A:72:GLN:HG2	1:A:75:ARG:HH21	1.78	0.48
5:E:36:ARG:NH1	5:E:65:TYR:OH	2.45	0.48
1:A:154:GLU:OE1	5:E:102:ARG:NH1	2.47	0.48
5:E:38:ASP:HA	5:E:39:PRO:HD2	1.65	0.48
1:A:268:LYS:HA	1:A:269:PRO:HD2	1.69	0.47
2:B:55:SER:OG	2:B:56:PHE:N	2.48	0.47
4:D:186:ASP:N	4:D:186:ASP:OD1	2.48	0.46
4:D:82:PRO:HA	4:D:114:VAL:HB	1.98	0.46
2:B:48:LYS:O	2:B:68:THR:OG1	2.31	0.46
4:D:138:VAL:HG11	5:E:146:VAL:HG21	1.98	0.45
1:A:223:ASP:OD1	1:A:223:ASP:N	2.35	0.45
2:B:64:LEU:HA	2:B:64:LEU:HD12	1.80	0.45
5:E:209:HIS:ND1	6:E:248:GOL:H31	2.33	0.44
5:E:177:GLN:HE21	5:E:177:GLN:HB2	1.56	0.44
2:B:96:ASP:HB3	2:B:99:MET:HB2	1.98	0.44
4:D:171:MET:HB3	4:D:173:PHE:HB2	1.99	0.44
4:D:120:ASN:HA	4:D:121:PRO:HD2	1.53	0.44
4:D:32:PHE:HD1	4:D:92:VAL:HG22	1.83	0.44
1:A:47:PRO:O	1:A:48:ARG:NH1	2.45	0.43
1:A:156:LEU:HA	1:A:156:LEU:HD23	1.77	0.43
1:A:195:SER:O	1:A:197:HIS:N	2.51	0.43
1:A:196:ASP:N	1:A:196:ASP:OD1	2.50	0.43
4:D:6:ASN:HA	4:D:6:ASN:HD22	1.64	0.42
4:D:153:SER:OG	4:D:158:VAL:O	2.37	0.42
5:E:29:HIS:ND1	5:E:94:SER:OG	2.48	0.42
4:D:201:ASP:OD2	4:D:201:ASP:N	2.53	0.41
4:D:154:LYS:HE3	4:D:154:LYS:HB3	1.89	0.41
2:B:1:ILE:HA	2:B:1:ILE:HD12	1.61	0.41
5:E:231:LYS:HA	5:E:232:PRO:HD3	1.87	0.41
5:E:148:LEU:HA	5:E:148:LEU:HD23	1.80	0.41
1:A:48:ARG:HA	1:A:48:ARG:HD3	1.82	0.40
5:E:59:GLU:H	5:E:59:GLU:HG3	1.69	0.40
2:B:71:THR:HA	2:B:72:PRO:HD2	1.88	0.40
5:E:27:MET:HE3	5:E:27:MET:HB3	1.97	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	273/275 (99%)	265 (97%)	6 (2%)	2 (1%)	30	62
2	B	99/100 (99%)	98 (99%)	1 (1%)	0	100	100
3	C	7/9 (78%)	6 (86%)	1 (14%)	0	100	100
4	D	198/200 (99%)	182 (92%)	15 (8%)	1 (0%)	38	70
5	E	243/245 (99%)	230 (95%)	11 (4%)	2 (1%)	27	58
All	All	820/829 (99%)	781 (95%)	34 (4%)	5 (1%)	33	66

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	196	ASP
4	D	6	ASN
5	E	39	PRO
5	E	183	PRO
1	A	269	PRO

### 5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	231/231 (100%)	211 (91%)	20 (9%)	15	33
2	B	96/95 (101%)	82 (85%)	14 (15%)	5	11
3	C	7/7 (100%)	4 (57%)	3 (43%)	0	0
4	D	178/178 (100%)	149 (84%)	29 (16%)	3	9
5	E	209/209 (100%)	178 (85%)	31 (15%)	4	11
All	All	721/720 (100%)	624 (86%)	97 (14%)	6	13

All (97) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	SER
1	A	12	VAL
1	A	32	GLN
1	A	35	ARG
1	A	45	MET
1	A	52	ILE
1	A	66	LYS
1	A	72	GLN
1	A	89	GLU
1	A	98	MET
1	A	113	TYR
1	A	115	GLN
1	A	182	THR
1	A	223	ASP
1	A	227	ASP
1	A	247	VAL
1	A	249	VAL
1	A	256	ARG
1	A	262	GLN
1	A	272	LEU
2	B	0	MET
2	B	1	ILE
2	B	6	LYS
2	B	34	ASP
2	B	44	GLU
2	B	50	GLU
2	B	64	LEU
2	B	68	THR
2	B	70	PHE
2	B	75	LYS
2	B	85	VAL
2	B	91	LYS
2	B	98	ASP
2	B	99	MET
3	C	1	LEU
3	C	6	VAL
3	C	7	ASN
4	D	1	LYS
4	D	6	ASN
4	D	23	THR
4	D	31	SER
4	D	39	SER

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Mol	Chain	Res	Type
4	D	41	LYS
4	D	51	SER
4	D	63	THR
4	D	76	LEU
4	D	87	THR
4	D	90	CYS
4	D	100	SER
4	D	103	LYS
4	D	110	THR
4	D	120	ASN
4	D	127	GLN
4	D	128	LEU
4	D	130	ASP
4	D	135	ASP
4	D	137	SER
4	D	145	ASP
4	D	151	SER
4	D	158	VAL
4	D	162	ASP
4	D	170	SER
4	D	183	ASN
4	D	186	ASP
4	D	195	ASN
4	D	197	ILE
5	E	7	THR
5	E	25	GLN
5	E	27	MET
5	E	31	TYR
5	E	55	THR
5	E	59	GLU
5	E	67	VAL
5	E	70	SER
5	E	72	THR
5	E	81	SER
5	E	88	SER
5	E	105	GLU
5	E	113	ARG
5	E	119	LEU
5	E	120	LYS
5	E	133	SER
5	E	137	ILE
5	E	138	SER

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Mol	Chain	Res	Type
5	E	145	LEU
5	E	146	VAL
5	E	163	VAL
5	E	166	LYS
5	E	170	SER
5	E	172	VAL
5	E	177	GLN
5	E	187	ASP
5	E	192	LEU
5	E	195	ARG
5	E	211	ARG
5	E	226	THR
5	E	246	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	GLN
4	D	5	GLN
4	D	6	ASN
4	D	105	GLN
4	D	111	GLN
4	D	119	GLN
4	D	127	GLN
5	E	177	GLN

### 5.3.3 RNA ⓘ

There are no RNA chains in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

8 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
6	GOL	A	276	-	5,5,5	0.51	0	5,5,5	0.75	0
6	GOL	A	277	-	5,5,5	0.48	0	5,5,5	0.73	0
6	GOL	A	278	-	5,5,5	0.44	0	5,5,5	0.90	0
6	GOL	A	279	-	5,5,5	0.32	0	5,5,5	0.54	0
6	GOL	A	280	-	5,5,5	0.41	0	5,5,5	0.40	0
7	SO4	B	100	-	4,4,4	0.59	0	6,6,6	0.52	0
6	GOL	E	247	-	5,5,5	0.32	0	5,5,5	0.90	0
6	GOL	E	248	-	5,5,5	0.47	0	5,5,5	0.58	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	GOL	A	276	-	-	0/4/4/4	0/0/0/0
6	GOL	A	277	-	-	0/4/4/4	0/0/0/0
6	GOL	A	278	-	-	0/4/4/4	0/0/0/0
6	GOL	A	279	-	-	0/4/4/4	0/0/0/0
6	GOL	A	280	-	-	0/4/4/4	0/0/0/0
7	SO4	B	100	-	-	0/0/0/0	0/0/0/0
6	GOL	E	247	-	-	0/4/4/4	0/0/0/0
6	GOL	E	248	-	-	0/4/4/4	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.



## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	275/275 (100%)	0.95	42 (15%) 3 3	42, 55, 65, 71	0
2	B	100/100 (100%)	0.67	12 (12%) 5 5	46, 54, 65, 83	0
3	C	9/9 (100%)	0.34	0 100 100	51, 56, 63, 63	0
4	D	200/200 (100%)	1.09	42 (21%) 1 2	42, 54, 66, 72	0
5	E	245/245 (100%)	0.61	31 (12%) 4 4	43, 54, 66, 72	0
All	All	829/829 (100%)	0.84	127 (15%) 3 3	42, 54, 66, 83	0

All (127) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	D	186	ASP	6.7
4	D	184	LYS	6.6
1	A	197	HIS	6.4
4	D	185	SER	6.2
1	A	223	ASP	6.0
4	D	134	SER	5.9
4	D	135	ASP	5.9
4	D	133	SER	5.9
4	D	154	LYS	5.9
4	D	171	MET	5.6
4	D	132	LYS	5.5
1	A	251	SER	5.3
4	D	153	SER	5.0
5	E	1	ASN	4.9
1	A	256	ARG	4.8
4	D	156	SER	4.7
1	A	257	TYR	4.6
1	A	195	SER	4.6
5	E	185	LEU	4.6
4	D	206	SER	4.5

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Mol	Chain	Res	Type	RSRZ
1	A	248	VAL	4.5
1	A	191	HIS	4.2
1	A	220	ASP	4.2
4	D	155	ASP	4.2
1	A	253	GLN	4.1
1	A	192	HIS	4.1
1	A	219	ARG	4.0
5	E	226	THR	4.0
4	D	197	ILE	4.0
1	A	218	GLN	3.9
1	A	221	GLY	3.9
1	A	226	GLN	3.9
1	A	272	LEU	3.9
1	A	275	GLU	3.8
5	E	41	MET	3.8
1	A	273	ARG	3.8
4	D	129	ARG	3.7
1	A	194	VAL	3.7
5	E	221	GLU	3.7
4	D	183	ASN	3.6
5	E	33	SER	3.6
1	A	196	ASP	3.5
1	A	58	GLU	3.5
1	A	254	GLU	3.5
1	A	255	GLN	3.5
1	A	222	GLU	3.5
5	E	184	ALA	3.5
4	D	187	PHE	3.4
1	A	199	ALA	3.4
5	E	166	LYS	3.4
4	D	172	ASP	3.4
5	E	2	ALA	3.4
4	D	169	ARG	3.3
1	A	250	PRO	3.3
1	A	274	TRP	3.3
1	A	252	GLY	3.2
4	D	200	GLU	3.2
5	E	186	ASN	3.2
1	A	224	GLN	3.1
5	E	245	ALA	3.0
4	D	170	SER	3.0
2	B	16	GLU	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	200	THR	3.0
4	D	191	ASN	3.0
1	A	1	GLY	2.9
2	B	80	CYS	2.9
5	E	246	ASP	2.8
5	E	222	ASN	2.8
4	D	145	ASP	2.8
5	E	183	PRO	2.8
4	D	205	PRO	2.7
2	B	99	MET	2.7
4	D	188	ALA	2.7
5	E	169	HIS	2.7
2	B	25	CYS	2.7
4	D	151	SER	2.6
5	E	92	CYS	2.6
1	A	227	ASP	2.6
4	D	157	ASP	2.6
2	B	58	LYS	2.6
4	D	195	ASN	2.6
5	E	117	GLU	2.6
4	D	74	SER	2.5
5	E	139	HIS	2.5
5	E	228	ASP	2.5
2	B	0	MET	2.5
5	E	207	ARG	2.5
5	E	134	GLU	2.4
4	D	7	SER	2.4
5	E	39	PRO	2.4
4	D	204	PHE	2.4
4	D	194	ASN	2.4
2	B	23	LEU	2.4
4	D	152	GLN	2.4
5	E	32	MET	2.4
2	B	75	LYS	2.4
1	A	258	THR	2.4
4	D	159	TYR	2.3
1	A	94	THR	2.3
5	E	138	SER	2.3
5	E	201	THR	2.3
4	D	201	ASP	2.3
5	E	90	TYR	2.3
4	D	147	GLN	2.3

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Mol	Chain	Res	Type	RSRZ
4	D	203	PHE	2.3
1	A	268	LYS	2.3
2	B	15	ALA	2.2
5	E	187	ASP	2.2
1	A	270	LEU	2.2
4	D	162	ASP	2.2
4	D	139	CYS	2.2
1	A	260	HIS	2.2
4	D	8	GLY	2.2
1	A	225	THR	2.2
4	D	136	LYS	2.2
4	D	174	LYS	2.2
5	E	93	ALA	2.2
1	A	177	GLU	2.1
2	B	24	ASN	2.1
1	A	95	VAL	2.1
1	A	229	GLU	2.1
1	A	216	THR	2.1
5	E	34	TRP	2.1
2	B	17	ASN	2.1
5	E	224	GLU	2.1
2	B	22	PHE	2.0
5	E	40	GLY	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

## 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
6	GOL	E	248	6/6	0.39	-	63,64,64,65	6
7	SO4	B	100	5/5	0.28	-	73,73,74,74	0
6	GOL	A	276	6/6	0.25	-	58,59,61,61	0
6	GOL	A	277	6/6	0.29	-	73,74,74,75	0
6	GOL	A	280	6/6	0.13	-	80,80,80,81	0
6	GOL	A	279	6/6	0.46	-	77,79,80,80	0
6	GOL	E	247	6/6	0.25	-	62,64,65,65	0
6	GOL	A	278	6/6	0.16	-	62,62,63,64	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.